

```
37 39 40
              41
ring nodes :
              5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 29 30 31 32 33 34 35 36
   26 27 28
chain bonds:
   6-39 17-37 23-26 24-25 39-40 39-41
ring bonds :
   1-2 1-5 1-23 2-3 3-4 4-5 4-6 6-7 7-8 7-9 8-11 9-10 10-11
                                                                  11-24 12-13 12-16
   12-23 13-14 14-15 15-16 15-17 17-18 18-19 18-20 19-22 20-21 21-22 22-24 25-32
                                                     33-34
   25-36 26-27 26-31 27-28 28-29 29-30 30-31 32-33
                                                            34-35
                                                                  35-36
exact/norm bonds :
   1-2 1-5 1-23 2-3 3-4 4-5 4-6 6-7 7-8 7-9 8-11 9-10 10-11
                                                                  11-24 12-13 12-16
   12-23 13-14 14-15 15-16 15-17 17-18 17-37 18-19 18-20 19-22 20-21 21-22 22-24
exact bonds:
   6-39 23-26 24-25 39-40 39-41
normalized bonds:
   25-32 25-36 26-27 26-31 27-28 28-29 29-30 30-31 32-33 33-34 34-35 35-36
```

G1:C,0

Match level:
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom 26:Atom 27:Atom 28:Atom 29:Atom 30:Atom 31:Atom 32:Atom 33:Atom 34:Atom 35:Atom 36:Atom 37:Atom 39:CLASS 40:CLASS 41:CLASS

Page 4

10/622,322

Structure attributes must be viewed using STN Express query preparation.

=> s l1

SAMPLE SEARCH INITIATED 10:11:19 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 10 TO ITERATE

100.0% PROCESSED 10 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS:

11 TO 389

PROJECTED ANSWERS:

1 TO 80

L2 1 SEA SSS SAM L1

=> s l1 sss full

FULL SEARCH INITIATED 10:11:26 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 154 TO ITERATE

100.0% PROCESSED 154 ITERATIONS

7 ANSWERS

SEARCH TIME: 00.00.01

L3 7 SEA SSS FUL L1

=> file caplus COST IN U.S. DOLLARS

SINCE FILE TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

155.42 155.63

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FILE COVERS 1907 - 31 Mar 2004 VOL 140 ISS 14 FILE LAST UPDATED: 30 Mar 2004 (20040330/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 13

L4

3 L3

=> d ibib abs hitstr tot

03/31/2004

Habte

```
L4 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN
ACCESSION NUMBER:
                         2000:742098 CAPLUS
DOCUMENT NUMBER:
                         133:296323
TITLE:
                         Process for modifying polypyrrolic macrocycles via
                         1,3-dipolar cycloadditions
                         MacAlpine, Jill; Dolphin, David; Sternberg, Ethan D.
INVENTOR (S):
PATENT ASSIGNEE(S):
                         THE University of British Columbia (UBC), Can.
SOURCE:
                         PCT Int. Appl., 41 pp.
                         CODEN: PIXXD2
DOCUMENT TYPE:
                         Patent
LANGUAGE:
                         English
FAMILY ACC. NUM. COUNT:
PATENT INFORMATION:
```

	NO.		DATE				CATI		ο.	DATE			
WO 200	0061585	A2	20001019							2000	0414		
			AT, AU,		BA,	BÐ,	BG,	BR,	BY,	CA,	ÇH,	CN,	CR,
			DM, DZ,										
	ID, IL,	IN, IS,	JP, KE,	KG,	KP,	KR,	KZ,	LC,	LK,	LR,	LS,	LT,	LU,
	LV, MA,	MD, MG,	MK, MN,	MW,	MX,	NO,	NZ,	PL,	PT,	RO,	RU,	ŞD,	ŞΕ,
	SG, SI,	SK, SL,	TJ, TM,	TR,	TT,	TZ,	UA,	UG,	US,	UΖ,	VN,	YU,	ZA,
	ZW, AM,	AZ, BY,	KG, KZ,	MD,	RU,	ТJ,	TM						
RW	: GH, GM,	KE, LS,	MW, SD,	SL,	SZ,	ΤZ,	UG,	ZW,	ΑT,	BE,	CH,	CY,	DE,
	DK, ES,	FI, FR,	GB, GR,	ΙE,	ΙT,	LU,	MC,	NL,	PT,	SE,	BF,	ΒJ,	CF,
	CG, CI,	CM, GA,	GN, GW,	ML,	MR,	ΝE,	SN,	TD,	TG				
US 6620929 B1 20030916 US 2000~551160 20000414													
US 200	4019201	A1	20040129		US	5 20	03-63	2232	2	2003	0717		
PRIORITY AP	PLN. INFO).:		ţ	US 19	999-	1293	24P	P	1999	0414		
				τ	US 20	000-	55116	50	Al	2000	0414		
OTHER SOURC	E(S):	CAS	REACT 13	3:29	6323;	, MA	RPAT	133	: 296	323			
GI													

AB Methods of modifying polypyrrolic macrocycles by use of a 1,3-dipolar cycloaddn. to produce compds. for further derivatization to produce photosensitizing agents (I) [M = 2H, Ni(II), Cu(II), Zn, Sn, Ge, Si, Ga,

L4 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN (Continued) kN21, kN22, kN23, kN24) bis [propanedinitrilato]] (2-)], (SP-4-1)- (9CI) (CA INDEX NAME)

L4 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN (continued)
Al, Mn(III), Gd(III), In, Tc; R1-R6 = independently H, alkyl,
alkylcarboxylic acid or ester, =0, OH, NO2, NH2; taken together with
another ring, ring substituent or meso substituent = fused 5 - or
6-membered ring; Ph1-Ph4 independently = H, (un)substituted alkyl,
(un)substituted aryl, (un)substituted cycloalkyl, may be same or
different] of interest is described.

IT 301300-67-4P 301300-68-5P 301545-80-2P
RL: BAC (Biological activity or effector, except adverse); BSU
(Biological
study, unclassified); IMF (Industrial manufacture); SPN (Synthetic
preparation); THU (Therapeutic use); BIOL (Biological study); PREP
(Preparation); USES (Uses)
(process for modifying polypyrrolic macrocycles via 1,3-dipolar
cycloaddns.)

RN 301300-67-4 CAPLUS
CN Propanedinitrile, (15-oxo-10,20-diphenyl-21H,23H-porphin-5(15H)-ylidene)(9CI) (CA INDEX NAME)

RN 301300-68-5 CAPLUS
CN Propanedinitrile,
2,2'-(10,20-diphenyl-21H,23H-porphine-5,15-diylidene)bis(9CI) (CA INDEX NAME)

RN 301545-80-2 CAPLUS CN Zinc, [[2,2] (10,20 diphenyl-21H,23H-porphine-5,15-diylidene-

L4 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN ACCESSION NUMBER: 2000:359162 CAPLUS DOCUMENT NUMBER: 133:83372 TITLE: Synthesis and crystal structure of a cumulenic quinoidal porphyrin dimer with strong electronic absorption in the infrared AUTHOR (S): Blake, Iain M.; Rees, Leigh H.; Claridge, Tim D. W.; Anderson, Harry L. CORPORATE SOURCE: Department of Chemistry, Dyson Perrins Laboratory, University of Oxford, Oxford, OX1 3QY, UK Angewandte Chemie, International Edition (2000), SOURCE: 39(10), 1818-1821 CODEN: ACIEF5; ISSN: 1433-7851 PUBLISHER: Wiley-VCH Verlag GmbH DOCUMENT TYPE: Journal LANGUAGE: English GI

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT * AB The cumulenic porphyrin dimer I (R = 2,6-bis(tert-butyl)phenyl) was prepared in 34 % yield by Pd/Cu catalyzed Takahashi coupling of the dibromoporphyrin dimer II (R = 2,6-bis(tert-butyl)phenyl) with the malononitrile anion followed by oxidation with N-iodosuccinimide. It was also encouraging to attempt the preparation of a shorter quinoidal (R = 2,6-bis(tert-butyl)phenyl) by first preparing the meso-meso linked dibromo dimer III (R = 2,6-bis(tert-butyl)phenyl). IV was prepared in yield. The electronic absorption of I and IV illustrate the amazing increase in conjugation accompanying quinoidalization. I-2C5H5N was characterized by single crystal x-ray diffraction anal. The compds. were all characterized by NMR and UV-visible spectra. RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (preparation and crystal structure of) 278807-38-8 CAPLUS Zinc, [µ-[[2,2'-[1,2-ethenediylidenebis[10,20-bis[3,5-bis[1,1 dimethylethyl)phenyl]-21H,23H-porphine-15,5-diylidenekN21, kN22, kN23, kN24]]bis(propanedinitrilato)](4-))]di-, compd. with pyridine (1:2) (9CI) (CA INDEX NAME)

CRN 278807-33-3 CMF C104 H100 N12 Zn2 CCI CCS (Continued)

L4 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN

CM

CRN 110-86-1 CMF C5 H5 N

278807-33-3P

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (preparation, mol. structure and NMR of)

278807-33-3 CAPLUS

Zinc, $[\mu \cdot [[2,2] \cdot [1,2]]$ ethenediylidenebis [10,20]-bis [3,5]-bis [1,1]dimethylethyl) phenyl] -21H, 23H-porphine-15, 5-diylidenekN21, kN22, kN23, kN24] bis [propanedinitrilato]] (4-)]]di- (9CI) (CA INDEX NAME)

L4 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN

ACCESSION NUMBER: 1998:630177 CAPLUS

DOCUMENT NUMBER: 129:330586

TITLE:

Synthesis, Crystal Structure, and Electronic Structure of a 5,15-Dialkylideneporphyrin: A TCNQ/Porphyrin

Hybrid

Blake, Iain M.; Anderson, Harry L.; Beljonne, David; Bredas, Jean-Luc; Clegg, William AUTHOR (5):

CORPORATE SOURCE: Department of Chemistry, University of Oxford Dyson

Perrins Laboratory, Oxford, OX1 3QY, UK SOURCE: Journal of the American Chemical Society (1998),

120(41), 10764 10765

CODEN: JACSAT; ISSN: 0002-7863 PUBLISHER: American Chemical Society

DOCUMENT TYPE: Journal

LANGUAGE: English CASREACT 129:330586 OTHER SOURCE(S):

Bu-t

The authors describe the synthesis, crystal structure and electronic structure of a 5,15-dialkylideneporphyrin (I) - a TCNQ/porphyrin hybrid. 215313-79-4P ΙT

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (crystal structure; synthesis, crystal structure, and electronic structure of a 5,15-dialkylideneporphyrin a TCNQ/porphyrin hybrid) RN 215313-79-4 CAPLUS

CN Zinc, [[2,2'-[10,20-bis[3,5-bis(1,1-dimethylethyl)phenyl]-21H,23H-porphine-5,15-diylidene-kN21, kN22, kN23, kN24]bis[propanedini trilato]](2-)]-, (SP-4-1)-, compd. with pyridine (1:1) (9CI) (CA INDEX NAME)

Habte

L4 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN

(Continued)

REFERENCE COUNT: THIS

THERE ARE 31 CITED REFERENCES AVAILABLE FOR

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN (Continued)

CM 1

CRN 215313-77-2 CMF C54 H50 N8 Zn

CCI CCS

CM 2

CRN 110-86-1 CMF C5 H5 N

215313-77-2P

RL: SPN (Synthetic preparation); PREP (Preparation) (synthesis, crystal structure, and electronic structure of a

5,15-dialkylideneporphyrin- a TCNQ/porphyrin hybrid)

215313-77-2 CAPLUS

CN Zinc, [[2,2'-[10,20-bis[3,5-bis(1,1-dimethylethyl)phenyl]-21H,23H-porphine-5,15-diylidene-kN21, kN22, kN23, kN24) bis (propanedini trilato]](2-)]-, (SP-4-1)- (9CI) (CA INDEX NAME)

L4 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN (Continued)

REFERENCE COUNT:

30 THERE ARE 30 CITED REFERENCES AVAILABLE FOR

FORMAT

RECORD. ALL CITATIONS AVAILABLE IN THE RE

Page 2

NEWS WWW

CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

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FILE 'HOME' ENTERED AT 09:13:33 ON 31 MAR 2004

=> file reg
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 09:13:41 ON 31 MAR 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2004 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 30 MAR 2004 HIGHEST RN 669048-54-8 DICTIONARY FILE UPDATES: 30 MAR 2004 HIGHEST RN 669048-54-8

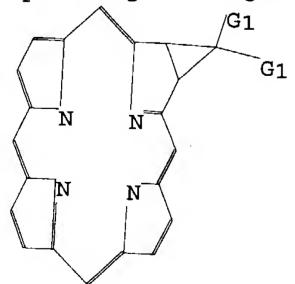
TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

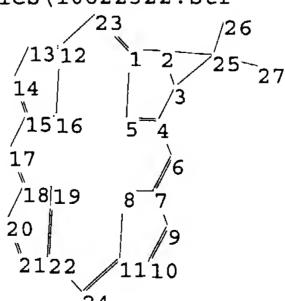
Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=> Uploading C:\Program Files\Stnexp\Queries\10622322.str





chain nodes : 26 27

Habte

03/31/2004

Page 3

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25

chain bonds: 25-26 25-27 ring bonds:

1-2 1-5 1-23 2-3 2-25 3-4 3-25 4-5 4-6 6-7 7-8 7-9 8-11 9-10 10-11 11-24 12-13 12-16 12-23 13-14 14-15 15-16 15-17 17-18 18-19 18-20 19-22

20-21 21-22 22-24

exact/norm bonds :

1-2 2-3 2-25 3-4 3-25 7-8 8-11 12-16 15-16 18-20 20-21 21-22 25-26 25-27

normalized bonds :

1-5 1-23 4-5 4-6 6-7 7-9 9-10 10-11 11-24 12-13 12-23 13-14 14-15

15-17 17-18 18-19 19-22 22-24

isolated ring systems :

containing 1 :

G1:CH2,CN

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom 26:CLASS

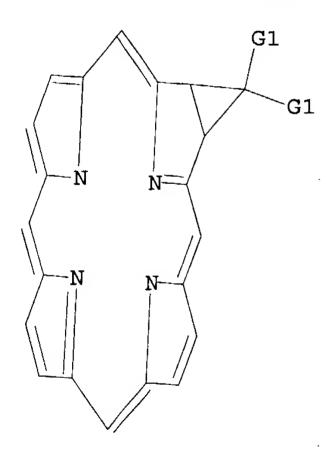
L1 STRUCTURE UPLOADED

=> d 11

L1 HAS NO ANSWERS

L1

STR



G1 CH2, CN

Structure attributes must be viewed using STN Express query preparation.

=> s 11

SAMPLE SEARCH INITIATED 09:14:04 FILE 'REGISTRY'

Habte

03/31/2004

10/622,322 Page 4

SAMPLE SCREEN SEARCH COMPLETED - 4 TO ITERATE

100.0% PROCESSED 4 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 4 TO 200

PROJECTED ITERATIONS: 4 TO 200
PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s l1 sss full FULL SEARCH INITIATED 09:14:10 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 123 TO ITERATE

100.0% PROCESSED 123 ITERATIONS 1 ANSWERS

SEARCH TIME: 00.00.01

L3 1 SEA SSS FUL L1

=> file caplus

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 155.42 155.63

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FILE COVERS 1907 - 31 Mar 2004 VOL 140 ISS 14 FILE LAST UPDATED: 30 Mar 2004 (20040330/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 13

L4 1 L3

=> d ibib abs hitstr

ACCESSION NUMBER:

DOCUMENT NUMBER:

Page 5

Process for modifying polypyrrolic macrocycles via TITLE: 1,3-dipolar cycloadditions MacAlpine, Jill; Dolphin, David; Sternberg, Ethan D. INVENTOR (S): THE University of British Columbia (UBC), Can. PATENT ASSIGNEE(S): SOURCE: PCT Int. Appl., 41 pp. CODEN: PIXXD2 DOCUMENT TYPE: Patent LANGUAGE: English FAMILY ACC. NUM. COUNT: PATENT INFORMATION: APPLICATION NO. DATE PATENT NO. KIND DATE 20001019 WO 2000-IB615 20000414 WO 2000061585 A2 WO 2000061585 A3 20010201 W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG US 6620929 B1 20030916 US 2000-551160 20000414 A1 20040129 US 2003-622322 20030717 US 2004019201 US 1999-129324P P 19990414 US 2000-551160 A1 20000414 PRIORITY APPLN. INFO.: CASREACT 133:296323; MARPAT 133:296323 OTHER SOURCE(S):

2000:742098 CAPLUS

L4 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2004 ACS on STN

133:296323

AB Methods of modifying polypyrrolic macrocycles by use of a 1,3-dipolar cycloaddn. to produce compds. for further derivatization to produce photosensitizing agents (I) [M = 2H, Ni(II), Cu(II), Zn, Sn, Ge, Si, Ga,

ANSWER 1 OF 1 CAPLUS COPYRIGHT 2004 ACS on STN (continued)
Al, Mn(III), Gd(III), In, Tc; R1 R6 = independently H, alkyl,
alkylcarboxylic acid or ester, =0, OH, NO2, NH2; taken together with
another ring, ring substituent or meso substituent = fused 5- or
6-membered ring; Ph1-Ph4 independently = H, (un)substituted alkyl,
(un)substituted aryl, (un)substituted cycloalkyl, may be same or
differentl of interest is described.

IT 301300-66-3P

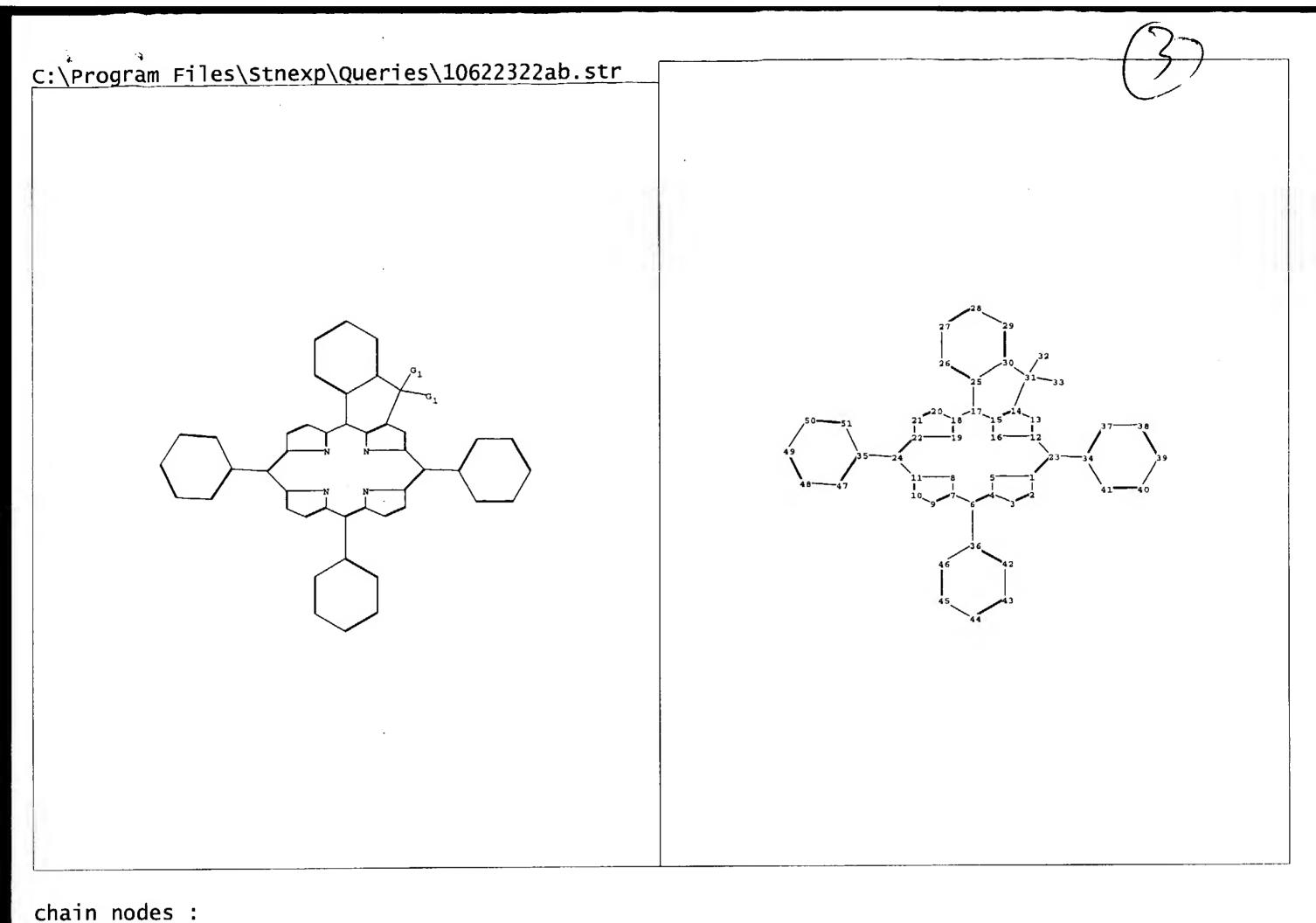
RL: BAC (Biological activity or effector, except adverse); BSU

(Biological study, unclassified); IMF (Industrial manufacture); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(process for modifying polypyrrolic macrocycles Via 1,3-dipolar cycloaddns.)

301300-66-3 CAPLUS

CN 1H,20H,22H-Cyclopropa[b]porphine-1,1-dicarbonitrile, 20,22-didehydro-1a,19a,21,23-tetrahydro-3,8,13,18-tetraphenyl- (9CI) (CA INDEX NAME)



```
15 16 17 18
40 41 42 43
                             35
                                 36
                                   37
                                       38
chain bonds :
   6-36 23-34 24-35 31-32 31-33
ring bonds:
   1-2 1-5 1-23 2-3 3-4 4-5 4-6 6-7 7-8 7-9 8-11 9-10 10-11
                                                                   11-24 12-13
                                                                                12-16
                      14-31 15-16 15-17 17-18 17-25
                                                       18-19
                                                                   19-22
                                                                          20-21
                                                                                21-22
                                                             18-20
   12-23 13-14 14-15
                                                                   35-47
                      26-27
                             27-28
                                  28-29 29-30 30-31
                                                                          35-51
                                                                                36-42
                                                      34-37 34-41
   22-24 25-26 25-30
                      39-40 40-41 42-43 43-44 44-45
                                                                                50-51
                                                             47 - 48
                                                                   48-49
                                                       45-46
                                                                          49-50
   36-46 37-38 38-39
exact/norm bonds :
                      3-4 4-5 4-6 6-7 7-8 7-9 8-11 9-10
                                                                                12-16
                                                             10-11
                                                                   11-24 12-13
   1-2 1-5 1-23 2-3
   12-23 13-14 14-15 14-31 15-16 15-17 17-18 17-25 18-19 18-20 19-22 20-21
                                                                               21-22
   22-24 30-31 31-32
                      31-33
exact bonds:
   6-36 23-34 24-35
normalized bonds:
   25-26 25-30 26-27 27-28 28-29 29-30 34-37 34-41 35-47 35-51 36-42 36-46 37-38
   38-39 39-40 40-41 42-43 43-44 44-45 45-46 47-48 48-49 49-50
                                                                   50-51
G1:CH2,CN,COOH
Match level:
   1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
                                                                              11:Atom
   12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom
                                                                             21:Atom
   22:Atom 23:Atom 24:Atom 25:Atom 26:Atom 27:Atom
                                                                              31:Atom
                                                    28:Atom 29:Atom
                                                                     30:Atom
           33:Atom 34:Atom 35:Atom 36:Atom 37:Atom
                                                                             41:Atom
                                                    38:Atom 39:Atom
                                                                     40:Atom
   32:Atom
                    44:Atom 45:Atom 46:Atom 47:Atom
                                                                              51:Atom
                                                                     50:Atom
                                                    48:Atom 49:Atom
   42:Atom
```

32 33

43:Atom

ring nodes:

	2					
						Q
		340				
			o'i			
			10			
			•			
1 0						
80			G.			
_						

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1

STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> s 11

SAMPLE SEARCH INITIATED 10:39:45 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 515 TO ITERATE

O ANSWERS

100.0% PROCESSED 515 ITERATIONS SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE** 8939 TO 11661

PROJECTED ITERATIONS:

PROJECTED ANSWERS:

0 TO

0 SEA SSS SAM L1 L2

=> s l1 sss full

FULL SEARCH INITIATED 10:39:52 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 10146 TO ITERATE

100.0% PROCESSED 10146 ITERATIONS

SEARCH TIME: 00.00.01

0 SEA SSS FUL L1

=> log y

L3

COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE ENTRY

155.42

SESSION 155.63

TOTAL

STN INTERNATIONAL LOGOFF AT 10:39:56 ON 31 MAR 2004

03/31/2004

Habte

Page 2

NEWS WWW

CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

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FILE 'HOME' ENTERED AT 10:38:55 ON 31 MAR 2004

=> file reg
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 10:39:09 ON 31 MAR 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2004 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 30 MAR 2004 HIGHEST RN 669048-54-8 DICTIONARY FILE UPDATES: 30 MAR 2004 HIGHEST RN 669048-54-8

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=>
Uploading C:\Program Files\Stnexp\Queries\10622322ab.str

Habte

L Number	Hits	Search Text	DB	Time stamp
1	648	540/145	USPAT; US-PGPUB	2004/03/31 14:13
3	276645	(carbonyl adj ylide)or synthesis	USPAT; US-PGPUB	2004/03/31 14:15
4	5021	photosensitizer	USPAT; US-PGPUB	2004/03/31 14:15
6	3449	tetraphenylporphyrin or diphenylporphyrin or TPP or DPP	USPAT	2004/03/31 14:16
7	31	540/145 and ((carbonyl adj ylide)or synthesis) and photosensitizer and (tetraphenylporphyrin or diphenylporphyrin or TPP or DPP)	USPAT	2004/03/31 14:17